



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 17, 2021 – 12:32 AM EDT

PDB ID : 1S7X
Title : Crystal structures of the murine class I major histocompatibility complex H-2Db in complex with LCMV-derived gp33 index peptide and three of its escape variants
Authors : Velloso, L.M.; Michaelsson, J.; Ljunggren, H.G.; Schneider, G.; Achour, A.
Deposited on : 2004-01-30
Resolution : 2.41 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.23.2
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

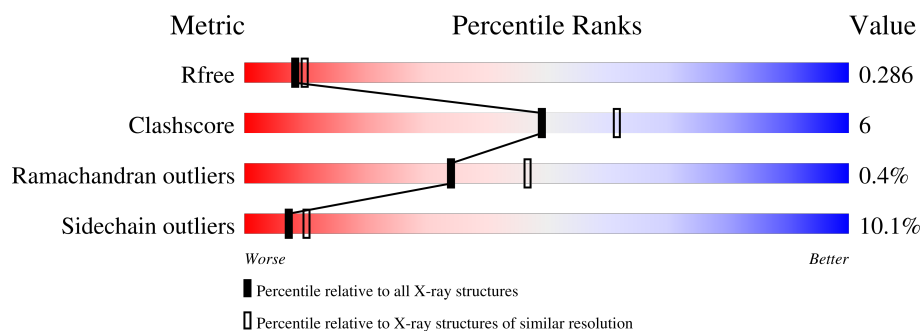
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION






The reported resolution of this entry is 2.41 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.




Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4647 (2.44-2.40)
Clashscore	141614	5161 (2.44-2.40)
Ramachandran outliers	138981	5073 (2.44-2.40)
Sidechain outliers	138945	5074 (2.44-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	338	
1	D	338	
1	G	338	
1	J	338	
2	B	99	
2	E	99	
2	H	99	

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Mol	Chain	Length	Quality of chain
2	K	99	 75% 22%
3	C	9	 89% 11%
3	F	9	 89% 11%
3	I	9	 67% 33%
3	L	9	 89% 11%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 13420 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called H-2 class I histocompatibility antigen, D-B alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	0	0
			2264	1430	400	425	9			
1	D	276	Total	C	N	O	S	0	0	0
			2264	1430	400	425	9			
1	G	276	Total	C	N	O	S	0	0	0
			2264	1430	400	425	9			
1	J	276	Total	C	N	O	S	0	0	0
			2264	1430	400	425	9			

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	0	0
			818	523	138	151	6			
2	E	99	Total	C	N	O	S	0	0	0
			818	523	138	151	6			
2	H	99	Total	C	N	O	S	0	0	0
			818	523	138	151	6			
2	K	99	Total	C	N	O	S	0	0	0
			818	523	138	151	6			

- Molecule 3 is a protein called Glycoprotein 9-residue peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	S	0	0	0
			72	48	11	12	1			
3	F	9	Total	C	N	O	S	0	0	0
			72	48	11	12	1			
3	I	9	Total	C	N	O	S	0	0	0
			72	48	11	12	1			
3	L	9	Total	C	N	O	S	0	0	0
			72	48	11	12	1			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	4	PHE	TYR	engineered mutation	UNP P07399
C	9	MET	CYS	SEE REMARK 999	UNP P07399
F	4	PHE	TYR	engineered mutation	UNP P07399
F	9	MET	CYS	SEE REMARK 999	UNP P07399
I	4	PHE	TYR	engineered mutation	UNP P07399
I	9	MET	CYS	SEE REMARK 999	UNP P07399
L	4	PHE	TYR	engineered mutation	UNP P07399
L	9	MET	CYS	SEE REMARK 999	UNP P07399

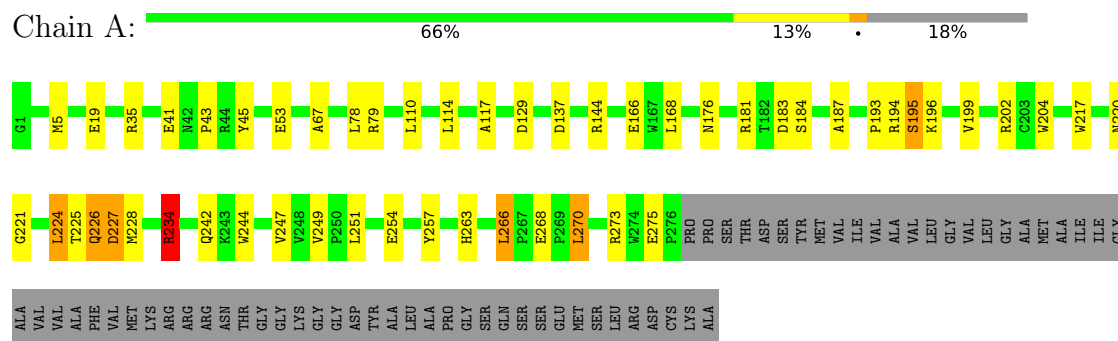
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	107	Total O 107 107	0	0
4	B	91	Total O 91 91	0	0
4	C	2	Total O 2 2	0	0
4	D	126	Total O 126 126	0	0
4	E	74	Total O 74 74	0	0
4	F	5	Total O 5 5	0	0
4	G	145	Total O 145 145	0	0
4	H	78	Total O 78 78	0	0
4	I	4	Total O 4 4	0	0
4	J	112	Total O 112 112	0	0
4	K	53	Total O 53 53	0	0
4	L	7	Total O 7 7	0	0

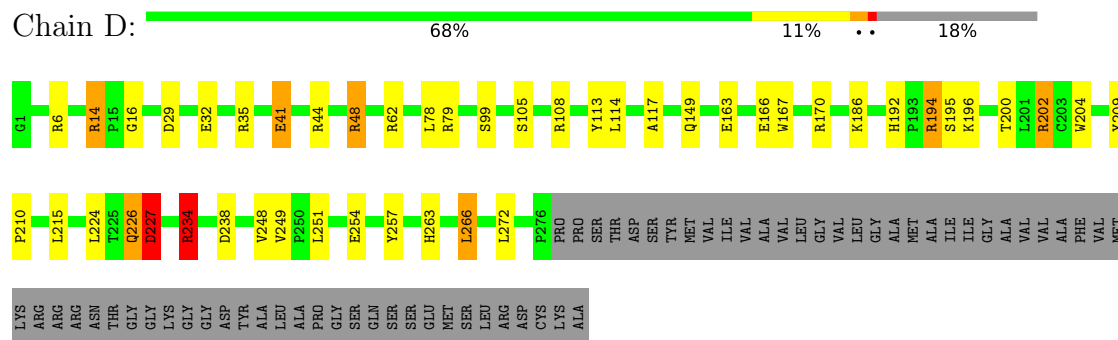
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

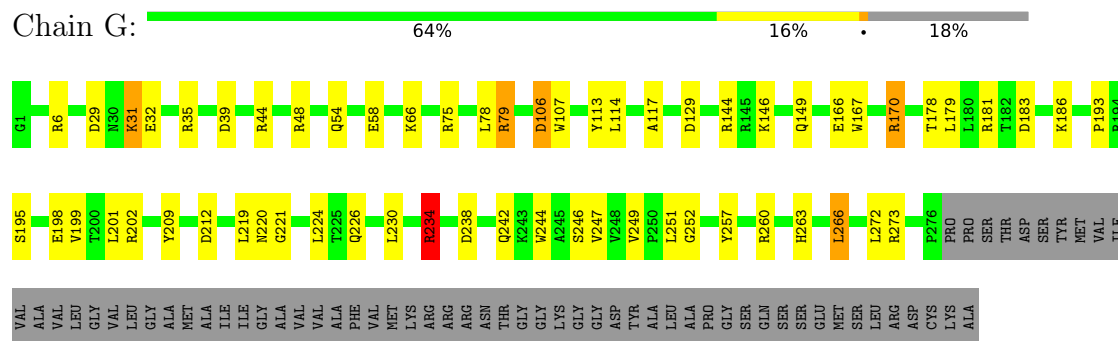
- Molecule 1: H-2 class I histocompatibility antigen, D-B alpha chain



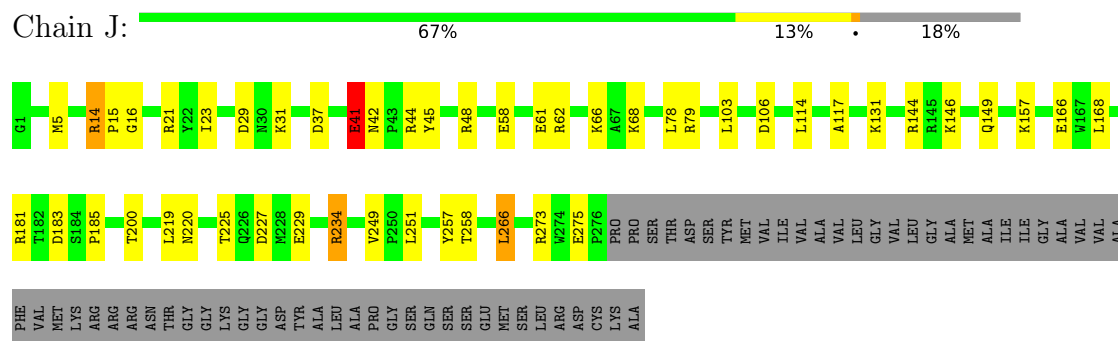
- Molecule 1: H-2 class I histocompatibility antigen, D-B alpha chain



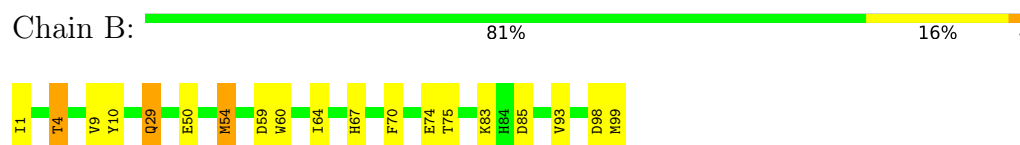
- Molecule 1: H-2 class I histocompatibility antigen, D-B alpha chain



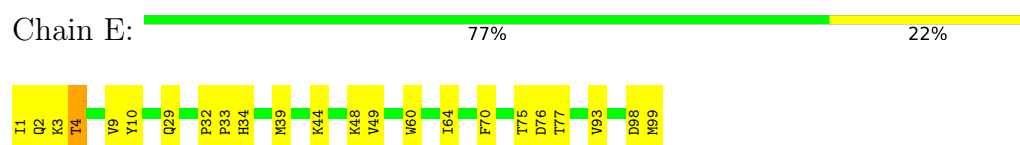
- Molecule 1: H-2 class I histocompatibility antigen, D-B alpha chain



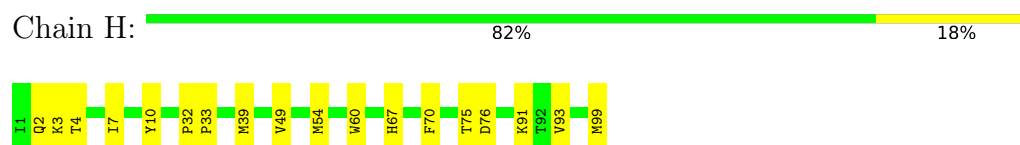
- Molecule 2: Beta-2-microglobulin



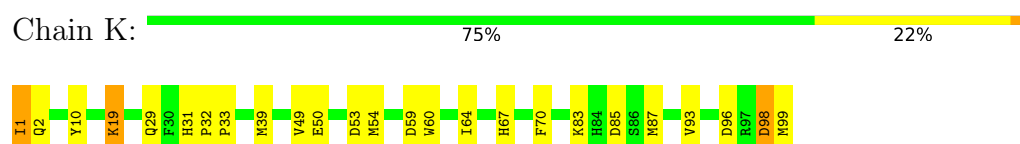
- Molecule 2: Beta-2-microglobulin



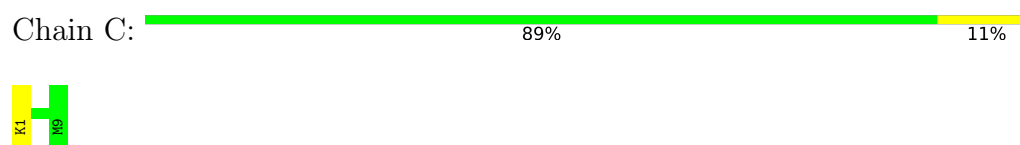
- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



- Molecule 3: Glycoprotein 9-residue peptide

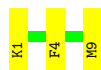


- Molecule 3: Glycoprotein 9-residue peptide

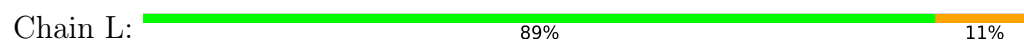




- Molecule 3: Glycoprotein 9-residue peptide



- Molecule 3: Glycoprotein 9-residue peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	91.90Å 123.32Å 100.20Å 90.00° 102.82° 90.00°	Depositor
Resolution (Å)	19.80 – 2.41 31.93 – 2.41	Depositor EDS
% Data completeness (in resolution range)	(Not available) (19.80-2.41) 98.9 (31.93-2.41)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.55 (at 2.42Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.203 , 0.254 0.245 , 0.286	Depositor DCC
R_{free} test set	2037 reflections (2.45%)	wwPDB-VP
Wilson B-factor (Å ²)	37.7	Xtriage
Anisotropy	0.082	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 57.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	13420	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.69% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.46	0/2331	0.74	6/3166 (0.2%)
1	D	0.51	0/2331	0.78	6/3166 (0.2%)
1	G	0.53	1/2331 (0.0%)	0.78	9/3166 (0.3%)
1	J	0.58	1/2331 (0.0%)	0.76	6/3166 (0.2%)
2	B	0.58	0/844	0.79	2/1146 (0.2%)
2	E	0.58	0/844	0.77	2/1146 (0.2%)
2	H	0.59	0/844	0.78	1/1146 (0.1%)
2	K	0.50	0/844	0.78	4/1146 (0.3%)
3	C	0.76	0/73	0.70	0/95
3	F	0.84	0/73	0.77	0/95
3	I	1.00	1/73 (1.4%)	0.87	0/95
3	L	0.72	0/73	0.67	0/95
All	All	0.54	3/12992 (0.0%)	0.77	36/17628 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	41	GLU	C-N	17.52	1.74	1.34
1	G	146	LYS	CE-NZ	6.06	1.64	1.49
3	I	9	MET	CG-SD	-5.05	1.68	1.81

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	41	GLU	CA-C-N	-9.36	96.60	117.20
1	J	41	GLU	O-C-N	7.65	134.94	122.70
1	A	227	ASP	CB-CG-OD2	6.43	124.09	118.30
1	A	129	ASP	CB-CG-OD2	6.42	124.07	118.30
1	G	29	ASP	CB-CG-OD2	6.18	123.86	118.30
2	K	98	ASP	CB-CG-OD2	6.11	123.80	118.30
1	D	234	ARG	NE-CZ-NH1	6.05	123.33	120.30
2	K	53	ASP	CB-CG-OD2	5.94	123.64	118.30
2	B	85	ASP	CB-CG-OD2	5.82	123.54	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	76	ASP	CB-CG-OD2	5.80	123.52	118.30
1	D	35	ARG	NE-CZ-NH2	-5.79	117.40	120.30
1	G	129	ASP	CB-CG-OD2	5.72	123.45	118.30
1	J	183	ASP	CB-CG-OD2	5.55	123.30	118.30
1	J	106	ASP	CB-CG-OD2	5.54	123.29	118.30
1	G	183	ASP	CB-CG-OD2	5.50	123.25	118.30
1	G	106	ASP	CB-CG-OD2	5.48	123.23	118.30
1	A	137	ASP	CB-CG-OD2	5.48	123.23	118.30
1	D	35	ARG	NE-CZ-NH1	5.47	123.03	120.30
1	G	146	LYS	CD-CE-NZ	5.43	124.18	111.70
2	E	76	ASP	CB-CG-OD2	5.32	123.09	118.30
1	A	35	ARG	NE-CZ-NH1	5.31	122.95	120.30
1	D	238	ASP	CB-CG-OD2	5.30	123.07	118.30
1	D	227	ASP	CB-CG-OD2	5.29	123.06	118.30
1	A	183	ASP	CB-CG-OD2	5.29	123.06	118.30
1	G	234	ARG	NE-CZ-NH1	5.28	122.94	120.30
2	E	98	ASP	CB-CG-OD2	5.28	123.05	118.30
1	D	29	ASP	CB-CG-OD2	5.26	123.03	118.30
2	B	98	ASP	CB-CG-OD2	5.23	123.00	118.30
2	K	59	ASP	CB-CG-OD2	5.22	123.00	118.30
1	A	234	ARG	NE-CZ-NH1	5.21	122.90	120.30
1	J	227	ASP	CB-CG-OD2	5.19	122.97	118.30
2	K	85	ASP	CB-CG-OD2	5.16	122.95	118.30
1	G	238	ASP	CB-CG-OD2	5.15	122.94	118.30
1	G	35	ARG	NE-CZ-NH1	5.05	122.83	120.30
1	G	212	ASP	CB-CG-OD2	5.04	122.83	118.30
1	J	29	ASP	CB-CG-OD2	5.03	122.83	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2264	0	2136	19	0
1	D	2264	0	2136	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	2264	0	2136	25	0
1	J	2264	0	2135	24	0
2	B	818	0	791	13	0
2	E	818	0	791	18	0
2	H	818	0	791	15	0
2	K	818	0	791	14	0
3	C	72	0	74	0	0
3	F	72	0	74	1	0
3	I	72	0	74	0	0
3	L	72	0	74	3	0
4	A	107	0	0	2	0
4	B	91	0	0	6	0
4	C	2	0	0	0	0
4	D	126	0	0	3	0
4	E	74	0	0	4	0
4	F	5	0	0	0	0
4	G	145	0	0	6	0
4	H	78	0	0	2	0
4	I	4	0	0	0	0
4	J	112	0	0	4	0
4	K	53	0	0	1	0
4	L	7	0	0	1	0
All	All	13420	0	12003	144	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (144) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:41:GLU:C	1:J:42:ASN:N	1.74	1.40
1:J:41:GLU:OE2	1:J:42:ASN:N	1.80	1.14
2:H:99:MET:C	4:H:102:HOH:O	1.88	1.12
2:B:99:MET:CG	4:B:180:HOH:O	2.01	1.07
1:D:14:ARG:HG2	1:D:14:ARG:HH11	0.92	1.05
2:E:99:MET:CG	4:E:105:HOH:O	2.09	1.00
1:D:14:ARG:HH11	1:D:14:ARG:CG	1.80	0.94
2:B:99:MET:C	4:B:101:HOH:O	2.07	0.91
1:D:14:ARG:HG2	1:D:14:ARG:NH1	1.73	0.89
2:E:99:MET:O	4:E:100:HOH:O	1.91	0.86
4:J:414:HOH:O	2:K:99:MET:CG	2.24	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:41:GLU:CD	1:D:41:GLU:H	1.83	0.80
1:J:41:GLU:CA	1:J:42:ASN:N	2.46	0.78
4:G:420:HOH:O	2:H:99:MET:CG	2.38	0.71
4:D:461:HOH:O	2:E:3:LYS:HE3	1.96	0.65
4:G:478:HOH:O	2:H:3:LYS:HE3	1.94	0.65
2:E:99:MET:C	4:E:100:HOH:O	2.36	0.63
4:G:478:HOH:O	2:H:3:LYS:CE	2.47	0.63
1:D:249:VAL:HG22	1:D:257:TYR:CZ	2.33	0.63
1:D:249:VAL:HG22	1:D:257:TYR:CE2	2.34	0.63
1:D:249:VAL:HG13	1:D:257:TYR:CE2	2.34	0.63
1:D:248:VAL:O	1:D:248:VAL:HG23	1.98	0.62
4:D:461:HOH:O	2:E:3:LYS:CE	2.48	0.61
2:K:39:MET:CE	2:K:49:VAL:HG13	2.31	0.60
1:A:144:ARG:HD2	4:A:342:HOH:O	2.01	0.60
1:G:32:GLU:OE2	1:G:48:ARG:HD2	2.02	0.60
1:J:21:ARG:HE	1:J:23:ILE:HD11	1.67	0.59
2:B:99:MET:O	4:B:101:HOH:O	2.14	0.59
1:D:117:ALA:HB2	2:E:60:TRP:CE2	2.38	0.59
2:K:1:ILE:N	4:K:121:HOH:O	2.34	0.59
1:A:234:ARG:HD3	2:B:10:TYR:CE2	2.37	0.59
1:D:32:GLU:OE2	1:D:48:ARG:HD2	2.03	0.59
1:G:201:LEU:HD12	1:G:249:VAL:HG21	1.84	0.58
1:J:185:PRO:HD2	1:J:266:LEU:HD13	1.88	0.56
1:J:144:ARG:HD2	4:J:429:HOH:O	2.06	0.55
1:J:234:ARG:HD3	2:K:10:TYR:CE2	2.41	0.55
1:A:194:ARG:O	1:A:195:SER:C	2.45	0.55
2:K:96:ASP:O	2:K:98:ASP:O	2.25	0.55
1:G:219:LEU:HG	1:G:220:ASN:HD22	1.72	0.55
4:G:420:HOH:O	2:H:99:MET:CB	2.54	0.54
2:H:39:MET:CE	2:H:49:VAL:HG13	2.38	0.54
1:G:224:LEU:HD23	1:G:247:VAL:HG21	1.90	0.54
1:J:41:GLU:CB	1:J:42:ASN:N	2.70	0.54
2:K:39:MET:HE2	2:K:49:VAL:HG13	1.90	0.53
1:G:117:ALA:HB2	2:H:60:TRP:CE2	2.46	0.51
1:G:193:PRO:HA	1:G:199:VAL:HG12	1.92	0.51
1:J:14:ARG:HG2	1:J:15:PRO:HD2	1.93	0.51
2:H:54:MET:HB2	4:H:101:HOH:O	2.09	0.51
1:A:249:VAL:HG22	1:A:257:TYR:CZ	2.45	0.51
1:D:194:ARG:HG2	4:D:353:HOH:O	2.11	0.51
2:K:98:ASP:C	2:K:99:MET:CG	2.79	0.51
2:H:39:MET:HE2	2:H:49:VAL:HG13	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:ARG:HD2	1:A:242:GLN:HB2	1.94	0.50
2:K:98:ASP:O	2:K:99:MET:CG	2.60	0.50
1:A:263:HIS:HB3	1:A:266:LEU:HD22	1.94	0.49
1:J:41:GLU:C	1:J:42:ASN:CA	2.75	0.49
1:D:234:ARG:HD3	2:E:10:TYR:CE2	2.47	0.49
1:A:224:LEU:HD23	1:A:247:VAL:HG21	1.95	0.49
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.47	0.48
2:H:7:ILE:HD12	2:H:91:LYS:HE2	1.95	0.48
1:G:202:ARG:HD2	1:G:244:TRP:CD2	2.49	0.48
3:L:1:LYS:NZ	4:L:731:HOH:O	2.46	0.48
1:J:21:ARG:NE	1:J:23:ILE:HD11	2.27	0.48
1:A:217:TRP:HB2	1:A:228:MET:CE	2.43	0.48
1:G:234:ARG:HD3	2:H:10:TYR:CE2	2.48	0.48
1:A:202:ARG:HD2	1:A:244:TRP:CD2	2.49	0.48
1:J:103:LEU:HD11	1:J:168:LEU:HD23	1.95	0.48
2:K:31:HIS:CD2	2:K:32:PRO:HA	2.49	0.48
2:B:29:GLN:NE2	2:B:59:ASP:OD2	2.47	0.47
1:D:263:HIS:HB3	1:D:266:LEU:HD22	1.96	0.47
2:B:9:VAL:HG23	2:B:93:VAL:HG22	1.96	0.47
1:D:234:ARG:HD3	2:E:10:TYR:CZ	2.49	0.47
1:G:144:ARG:NH2	4:G:398:HOH:O	2.47	0.47
1:J:117:ALA:HB2	2:K:60:TRP:CE2	2.49	0.47
1:G:249:VAL:HG22	1:G:257:TYR:CE1	2.50	0.47
1:D:14:ARG:NH2	2:E:34:HIS:HB2	2.30	0.46
2:E:39:MET:CE	2:E:49:VAL:HG13	2.45	0.46
1:D:14:ARG:CG	1:D:14:ARG:NH1	2.49	0.46
1:A:187:ALA:HA	1:A:204:TRP:O	2.15	0.46
1:G:251:LEU:HD23	1:G:252:GLY:N	2.31	0.46
1:G:201:LEU:O	1:G:246:SER:HA	2.16	0.46
1:J:5:MET:HB2	1:J:168:LEU:HD13	1.98	0.46
1:A:226:GLN:HA	1:A:227:ASP:HA	1.72	0.46
2:E:32:PRO:HB2	2:E:33:PRO:HD2	1.98	0.46
1:G:249:VAL:HG22	1:G:257:TYR:CZ	2.50	0.46
1:D:14:ARG:HH21	2:E:34:HIS:HB2	1.80	0.45
1:D:202:ARG:HG2	1:D:204:TRP:NE1	2.32	0.45
1:G:249:VAL:HG13	1:G:257:TYR:CE2	2.52	0.45
1:D:14:ARG:HH21	2:E:34:HIS:CG	2.34	0.45
2:H:39:MET:HE1	2:H:67:HIS:C	2.36	0.45
2:E:9:VAL:HG23	2:E:93:VAL:HG22	1.98	0.45
1:J:66:LYS:HZ2	3:L:1:LYS:HE2	1.81	0.45
1:J:131:LYS:NZ	1:J:157:LYS:HE3	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:178:THR:HG22	1:G:178:THR:O	2.18	0.44
1:A:5:MET:HB2	1:A:168:LEU:HD13	2.00	0.44
1:G:79:ARG:NE	4:G:388:HOH:O	2.51	0.43
1:D:14:ARG:NH2	2:E:34:HIS:CB	2.81	0.43
1:G:263:HIS:HB3	1:G:266:LEU:HD22	2.00	0.43
2:B:4:THR:HG22	4:B:104:HOH:O	2.19	0.43
1:J:41:GLU:HB2	1:J:42:ASN:N	2.33	0.43
1:G:234:ARG:HD2	1:G:242:GLN:HB2	2.01	0.43
1:G:224:LEU:CD2	1:G:247:VAL:HG21	2.49	0.43
2:E:4:THR:HG22	4:E:112:HOH:O	2.18	0.43
1:G:31:LYS:HG3	1:G:209:TYR:OH	2.18	0.43
1:J:249:VAL:HG22	1:J:257:TYR:CE1	2.54	0.43
1:J:66:LYS:HZ2	3:L:1:LYS:CE	2.33	0.42
2:B:54:MET:HB2	4:B:110:HOH:O	2.19	0.42
1:J:14:ARG:HH11	1:J:16:GLY:H	1.68	0.42
2:K:50:GLU:HB2	2:K:67:HIS:CE1	2.54	0.42
1:A:45:TYR:CE2	1:A:67:ALA:HB2	2.55	0.42
1:D:99:SER:HA	1:D:113:TYR:O	2.19	0.42
2:B:50:GLU:HB2	2:B:67:HIS:CE1	2.55	0.41
1:G:230:LEU:C	1:G:230:LEU:HD12	2.41	0.41
1:D:192:HIS:HB2	1:D:200:THR:HB	2.02	0.41
1:D:226:GLN:O	1:D:227:ASP:C	2.58	0.41
1:G:6:ARG:NH2	1:G:113:TYR:CE1	2.88	0.41
1:J:21:ARG:NH2	1:J:37:ASP:OD2	2.53	0.41
1:A:234:ARG:HD3	2:B:10:TYR:CD2	2.56	0.41
1:D:14:ARG:HH21	2:E:34:HIS:CB	2.33	0.41
1:G:167:TRP:HZ3	1:G:170:ARG:HE	1.68	0.41
2:H:7:ILE:HD12	2:H:91:LYS:CE	2.50	0.41
1:D:209:TYR:CD1	1:D:210:PRO:HA	2.56	0.41
1:G:106:ASP:O	1:G:107:TRP:HB2	2.20	0.41
1:A:270:LEU:HD12	4:A:377:HOH:O	2.20	0.41
1:D:251:LEU:HD23	1:D:254:GLU:OE1	2.21	0.41
2:H:54:MET:O	2:H:54:MET:HG3	2.20	0.41
1:D:163:GLU:OE1	3:F:1:LYS:NZ	2.53	0.41
1:G:201:LEU:CD1	1:G:249:VAL:HG21	2.50	0.41
1:A:193:PRO:HA	1:A:199:VAL:HG12	2.02	0.41
2:B:9:VAL:CG2	2:B:93:VAL:HG22	2.51	0.41
2:B:99:MET:CB	4:B:180:HOH:O	2.54	0.41
1:J:41:GLU:HB2	4:J:450:HOH:O	2.20	0.41
1:J:62:ARG:HG3	4:J:384:HOH:O	2.20	0.41
2:K:19:LYS:HB3	2:K:19:LYS:NZ	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:6:ARG:NH2	1:D:113:TYR:CE1	2.89	0.40
1:D:32:GLU:OE2	1:D:48:ARG:CD	2.68	0.40
2:E:48:LYS:HB3	2:E:48:LYS:NZ	2.35	0.40
1:J:41:GLU:OE2	1:J:42:ASN:CA	2.64	0.40
2:K:39:MET:HE3	2:K:49:VAL:HG13	2.03	0.40
1:A:41:GLU:O	1:A:43:PRO:HD3	2.22	0.40
1:A:249:VAL:HG22	1:A:257:TYR:CE1	2.57	0.40
1:D:167:TRP:CE3	1:D:170:ARG:HD3	2.56	0.40
2:H:32:PRO:HB2	2:H:33:PRO:HD2	2.02	0.40
2:K:32:PRO:HB2	2:K:33:PRO:HD2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	274/338 (81%)	258 (94%)	14 (5%)	2 (1%)	22	31
1	D	274/338 (81%)	263 (96%)	9 (3%)	2 (1%)	22	31
1	G	274/338 (81%)	253 (92%)	20 (7%)	1 (0%)	34	47
1	J	274/338 (81%)	260 (95%)	13 (5%)	1 (0%)	34	47
2	B	97/99 (98%)	94 (97%)	3 (3%)	0	100	100
2	E	97/99 (98%)	95 (98%)	2 (2%)	0	100	100
2	H	97/99 (98%)	95 (98%)	2 (2%)	0	100	100
2	K	97/99 (98%)	94 (97%)	3 (3%)	0	100	100
3	C	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
3	F	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
3	I	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
3	L	7/9 (78%)	6 (86%)	1 (14%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1512/1784 (85%)	1436 (95%)	70 (5%)	6 (0%)	34	47

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	227	ASP
1	G	221	GLY
1	D	16	GLY
1	A	195	SER
1	J	41	GLU
1	A	221	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	234/280 (84%)	211 (90%)	23 (10%)	8	11
1	D	234/280 (84%)	211 (90%)	23 (10%)	8	11
1	G	234/280 (84%)	210 (90%)	24 (10%)	7	10
1	J	234/280 (84%)	207 (88%)	27 (12%)	5	7
2	B	93/94 (99%)	84 (90%)	9 (10%)	8	11
2	E	93/94 (99%)	84 (90%)	9 (10%)	8	11
2	H	93/94 (99%)	88 (95%)	5 (5%)	22	34
2	K	93/94 (99%)	83 (89%)	10 (11%)	6	8
3	C	7/7 (100%)	6 (86%)	1 (14%)	3	3
3	F	7/7 (100%)	6 (86%)	1 (14%)	3	3
3	I	7/7 (100%)	5 (71%)	2 (29%)	0	0
3	L	7/7 (100%)	6 (86%)	1 (14%)	3	3
All	All	1336/1524 (88%)	1201 (90%)	135 (10%)	7	10

All (135) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	GLU
1	A	53	GLU
1	A	78	LEU
1	A	79	ARG
1	A	110	LEU
1	A	114	LEU
1	A	166	GLU
1	A	176	ASN
1	A	181	ARG
1	A	184	SER
1	A	196	LYS
1	A	220	ASN
1	A	224	LEU
1	A	225	THR
1	A	226	GLN
1	A	234	ARG
1	A	251	LEU
1	A	254	GLU
1	A	266	LEU
1	A	268	GLU
1	A	270	LEU
1	A	273	ARG
1	A	275	GLU
2	B	1	ILE
2	B	4	THR
2	B	29	GLN
2	B	54	MET
2	B	64	ILE
2	B	70	PHE
2	B	74	GLU
2	B	75	THR
2	B	83	LYS
3	C	1	LYS
1	D	14	ARG
1	D	41	GLU
1	D	44	ARG
1	D	48	ARG
1	D	62	ARG
1	D	78	LEU
1	D	79	ARG
1	D	105	SER
1	D	108	ARG
1	D	114	LEU

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Mol	Chain	Res	Type
1	D	149	GLN
1	D	166	GLU
1	D	186	LYS
1	D	194	ARG
1	D	195	SER
1	D	196	LYS
1	D	202	ARG
1	D	215	LEU
1	D	224	LEU
1	D	226	GLN
1	D	234	ARG
1	D	266	LEU
1	D	272	LEU
2	E	1	ILE
2	E	2	GLN
2	E	4	THR
2	E	29	GLN
2	E	44	LYS
2	E	64	ILE
2	E	70	PHE
2	E	75	THR
2	E	77	THR
3	F	1	LYS
1	G	31	LYS
1	G	39	ASP
1	G	44	ARG
1	G	54	GLN
1	G	58	GLU
1	G	66	LYS
1	G	75	ARG
1	G	78	LEU
1	G	79	ARG
1	G	114	LEU
1	G	149	GLN
1	G	166	GLU
1	G	170	ARG
1	G	179	LEU
1	G	181	ARG
1	G	186	LYS
1	G	195	SER
1	G	198	GLU
1	G	226	GLN

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Mol	Chain	Res	Type
1	G	234	ARG
1	G	260	ARG
1	G	266	LEU
1	G	272	LEU
1	G	273	ARG
2	H	2	GLN
2	H	4	THR
2	H	70	PHE
2	H	75	THR
2	H	93	VAL
3	I	1	LYS
3	I	4	PHE
1	J	14	ARG
1	J	31	LYS
1	J	41	GLU
1	J	44	ARG
1	J	45	TYR
1	J	48	ARG
1	J	58	GLU
1	J	61	GLU
1	J	68	LYS
1	J	78	LEU
1	J	79	ARG
1	J	114	LEU
1	J	146	LYS
1	J	149	GLN
1	J	166	GLU
1	J	181	ARG
1	J	200	THR
1	J	219	LEU
1	J	220	ASN
1	J	225	THR
1	J	229	GLU
1	J	234	ARG
1	J	251	LEU
1	J	258	THR
1	J	266	LEU
1	J	273	ARG
1	J	275	GLU
2	K	1	ILE
2	K	2	GLN
2	K	19	LYS

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Mol	Chain	Res	Type
2	K	29	GLN
2	K	54	MET
2	K	64	ILE
2	K	70	PHE
2	K	83	LYS
2	K	87	MET
2	K	93	VAL
3	L	1	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	97	GLN
1	A	192	HIS
1	A	226	GLN
2	B	38	GLN
3	C	5	ASN
1	D	54	GLN
1	D	97	GLN
2	E	67	HIS
3	F	5	ASN
1	G	54	GLN
1	G	97	GLN
1	G	220	ASN
2	H	67	HIS
3	I	5	ASN
1	J	30	ASN
1	J	72	GLN
1	J	97	GLN
1	J	192	HIS
1	J	220	ASN
2	K	31	HIS
2	K	67	HIS
3	L	5	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	J	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	J	41:GLU	C	42:ASN	N	1.74

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.